



# SEQUENCE LISTING

<110> University of British Columbia, et al.

<120> CXCR4 AGONIST TREATMENT OF HEMATOPOIETIC CELLS

<130> 80021-255

<140> US 09/835,107

<141> 2001-04-12

<150> CA 2,305,036

<151> 2000-04-12

<150> US 60/232,425

<151> 2000-09-14

<150> CA 2,335,109

<151> 2001-02-23

<160> 31

<170> PatentIn Ver. 2.0

<210> 1

<211> 67

<212> PRT

<213> Homo sapiens

<220>

<223> SDF-1 alpha

<220>

<221> MISC\_FEATURE

<222> (1)..(67)

<223> A pegylation moiety may be provided at any position on the sequence.

<400> 1

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
1 5 10 15

His Val Ala Arg Ala Asn Val Lys His Leu Lys Ile Leu Asn Thr Pro  
20 25 30

Asn Cys Ala Leu Gln Ile Val Ala Arg Leu Lys Asn Asn Asn Arg Gln  
35 40 45

Val Cys Ile Asp Pro Lys Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys  
50 55 60

Ala Leu Asn  
65

<210> 2

<211> 93

<212> PRT

<213> Homo sapiens

<220>

<223> SDF-1 Precursor, PBSF

<220>

<221> MISC\_FEATURE

<222> (1)..(93)

<223> A pegylation moiety may be provided at any position on the sequence.

<400> 2

Met Asn Ala Lys Val Val Val Val Leu Val Leu Val Leu Thr Ala Leu  
1 5 10 15

Cys Leu Ser Asp Gly Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys  
20 25 30

Arg Phe Phe Glu Ser His Val Ala Arg Ala Asn Val Lys His Leu Lys  
35 40 45

Ile Leu Asn Thr Pro Asn Cys Ala Leu Gln Ile Val Ala Arg Leu Lys  
50 55 60

Asn Asn Asn Arg Gln Val Cys Ile Asp Pro Lys Leu Lys Trp Ile Gln  
65 70 75 80

Glu Tyr Leu Glu Lys Ala Leu Asn Lys Arg Phe Lys Met  
85 90

<210> 3

<211> 93

<212> PRT

<213> Homo sapiens

<220>

<223> SDF-1 beta

<220>

<221> MISC\_FEATURE

<222> (1)..(93)

<223> A pegylation moiety may be provided at any position on the sequence.

<400> 3

Met Asn Ala Lys Val Val Val Val Leu Val Leu Val Leu Thr Ala Leu  
1 5 10 15

Cys Leu Ser Asp Gly Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys  
20 25 30

Arg Phe Phe Glu Ser His Val Ala Arg Ala Asn Val Lys His Leu Lys  
35 40 45

Ile Leu Asn Thr Pro Asn Cys Ala Leu Gln Ile Val Ala Arg Leu Lys  
50 55 60

Asn Asn Asn Arg Gln Val Cys Ile Asp Pro Lys Leu Lys Trp Ile Gln  
65 70 75 80

Glu Tyr Leu Glu Lys Ala Leu Asn Lys Arg Phe Lys Met  
85 90

<210> 4  
<211> 17  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory: SDF-1(1-17): or  
CTCE9902

<400> 4  
Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
1 5 10 15

His

<210> 5  
<211> 6  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory

<400> 5  
Arg Phe Phe Glu Ser His  
1 5

<210> 6  
<211> 9  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory

<400> 6  
Lys Pro Val Ser Leu Ser Tyr Arg Cys  
1 5

<210> 7  
<211> 9  
<212> PRT  
<213> Artificial Sequence

<220>  
<221> DISULFID  
<222> (9)  
<223> Disulphide linkage may form between two cys  
residues at position 9 of each of two monomers  
thereby forming a dimer.

<220>  
<223> Synthesised in Laboratory:  
SDF-1(1-9)2-C9/C9-cysteine dimer: or CTCE9901

<400> 7  
Lys Pro Val Ser Leu Ser Tyr Arg Cys  
1 5

<210> 8  
<211> 9  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory

<220>  
<221> BINDING  
<222> (9)  
<223> Linking Moiety (may be lysine with both the alpha and the epsilon amino groups of the lysine being associated with the covalent (amide) bond formation) may bind here allowing formation of a dimer.

<400> 8  
Lys Pro Val Ser Leu Ser Tyr Arg Cys  
1 5

<210> 9  
<211> 8  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory

<220>  
<221> BINDING  
<222> (8)  
<223> Linking Moiety (may be lysine with both the alpha and the epsilon amino groups of the lysine being associated with the covalent (amide) bond formation) may bind here allowing formation of a dimer.

<400> 9  
Lys Pro Val Ser Leu Ser Tyr Arg  
1 5

<210> 10  
<211> 30  
<212> PRT  
<213> Artificial Sequence

<220>  
<221> DOMAIN  
<222> (15)..(17)  
<223> spacer monomers (such as the illustrated glycine G's) may be used in variable numbers, such as 2, 3 or 4 glycines.

<220>  
<223> Synthesised in Laboratory:

SDF-1(1-14)-(G)3-SDF-1(55-67) acid

<400> 10  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
           1                  5                  10                  15

Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
                   20                  25                  30

<210> 11  
 <211> 31  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (16)..(19)  
 <223> spacer monomers (such as the illustrated glycine  
           G's) may be used in variable numbers, such as 2, 3  
           or 4 glycines.

<220>  
 <223> Synthesised in Laboratory:  
           SDF-1(1-14)-(G)4-SDF-1(55-67) acid: or CTCE0013

<400> 11  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
           1                  5                  10                  15

Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
                   20                  25                  30

<210> 12  
 <211> 30  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (15)..(17)  
 <223> spacer monomers (such as the illustrated glycine  
           G's) may be used in variable numbers, such as 2, 3  
           or 4 glycines.

<220>  
 <223> Synthesised in Laboratory:  
           SDF-1(1-14)-(G)3-SDF-1(55-67) amide

<220>  
 <221> MOD\_RES  
 <222> (30)  
 <223> AMIDATION

<400> 12  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
           1                  5                  10                  15

Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
                   20                  25                  30

<210> 13  
<211> 31  
<212> PRT  
<213> Artificial Sequence

<220>  
<221> DOMAIN  
<222> (15)..(18)  
<223> spacer monomers (such as the illustrated glycine  
G's) may be used in variable numbers, such as 2, 3  
or 4 glycines.

<220>  
<223> Synthesised in Laboratory:  
SDF-1(1-14)-(G)4-SDF-1(55-67) amide: or CTCE0017

<220>  
<221> MOD\_RES  
<222> (31)  
<223> AMIDATION

<400> 13  
Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15

Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
20 25 30

<210> 14  
<211> 33  
<212> PRT  
<213> Artificial Sequence

<220>  
<221> DOMAIN  
<222> (18)..(21)  
<223> spacer monomers (such as the illustrated glycine  
G's) may be used in variable numbers, such as 2, 3  
or 4 glycines.

<220>  
<223> Synthesised in Laboratory:  
SDF-1(1-17)-(G)3-SDF-1(55-67) acid

<400> 14  
Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
1 5 10 15

His Gly Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu  
20 25 30

Asn

<210> 15  
<211> 34  
<212> PRT

<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (18)..(21)

<223> spacer monomers (such as the illustrated glycine G's) may be used in variable numbers, such as 2, 3 or 4 glycines.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-17)-(G)4-SDF-1(55-67) acid

<400> 15

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
1 5 10 15

His Gly Gly Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala  
20 25 30

Leu Asn

<210> 16

<211> 33

<212> PRT

<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (18)..(20)

<223> spacer monomers (such as the illustrated glycine G's) may be used in variable numbers, such as 2, 3 or 4 glycines.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-17)-(G)3-SDF-1(55-67) amide

<220>

<221> MOD\_RES

<222> (33)

<223> AMIDATION

<400> 16

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
1 5 10 15

His Gly Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu  
20 25 30

Asn

<210> 17

<211> 34

<212> PRT

<213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (18)..(21)  
 <223> spacer monomers (such as the illustrated glycine  
 G's) may be used in variable numbers, such as 2, 3  
 or 4 glycines.

<220>  
 <223> Synthesised in Laboratory:  
 SDF-1(1-17)-(G)3-SDF-1(55-67) amide

<220>  
 <221> MOD\_RES  
 <222> (34)  
 <223> AMIDATION

<400> 17  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Glu Ser  
 1 5 10 15  
 His Gly Gly Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala  
 20 25 30  
 Leu Asn

<210> 18  
 <211> 31  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (15)..(18)  
 <223> spacer monomers (such as the illustrated glycine  
 G's) may be used in variable numbers, such as 2, 3  
 or 4 glycines.

<220>  
 <221> DOMAIN  
 <222> (24)..(28)  
 <223> Cyclized, for example glutamate (E) and lysine (K)  
 residues may be joined by side chain cyclization  
 using a lactam formation procedure.

<220>  
 <223> Synthesised in Laboratory:  
 SDF-1(1-14)-(G)4-SDF(55-67)-E24/K28-cyclic acid

<400> 18  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
 1 5 10 15  
 Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
 20 25 30

<210> 19  
 <211> 31  
 <212> PRT



<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (15)..(18)

<223> spacer monomers (such as the illustrated glycine G's) may be used in variable numbers, such as 2, 3 or 4 glycines.

<220>

<221> DOMAIN

<222> (20)..(24)

<223> Cyclized, for example glutamate (E) and lysine (K) residues may be joined by side chain cyclization using a lactam formation.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-14) - (G)4-SDF-1(55-67) -K20/E24-cyclic acid

<400> 19

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15

Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
20 25 30

<210> 20

<211> 31

<212> PRT

<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (15)..(18)

<223> spacer monomers (such as the illustrated glycine G's) may be used in variable numbers, such as 2, 3 or 4 glycines.

<220>

<221> DOMAIN

<222> (24)..(28)

<223> Cyclized, for example (E) and lysine (K) residues may be joined by side chain cyclization using a lactam formation procedure.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-14) - (G)4-SDF-1(55-67) -E24/K28-cyclic amide: or CTCE0022

<220>

<221> MOD\_RES

<222> (31)

<223> AMIDATION

<400> 20

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15

Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
                   20                                  25                                  30

<210> 21  
 <211> 31  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (15)..(18)  
 <223> spacer monomers (such as the illustrated glycine  
           G's) may be used in variable numbers, such as 2, 3  
           or 4 glycines.

<220>  
 <221> DOMAIN  
 <222> (20)..(24)  
 <223> Cyclized, for example glutamate (E) and lysine  
           (K) residues may be joined by side chain  
           cyclization using a lactam formation procedure.

<220>  
 <223> Synthesised in Laboratory:  
           SDF-1(1-14)-(G)4-SDF-1(55-67)-K20/E24-cyclic  
           amide: or CTCE0021

<220>  
 <221> MOD\_RES  
 <222> (31)  
 <223> AMIDATION

<400> 21  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
           1                                  5                                  10                                  15

Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
                   20                                  25                                  30

<210> 22  
 <211> 31  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (15)..(18)  
 <223> spacer monomers (such as the illustrated glycine  
           G's) may be used in variable numbers, such as 2, 3  
           or 4 glycines.

<220>  
 <221> DOMAIN  
 <222> (20)..(24)  
 <223> Internal cyclization of peptides of the invention  
           may be in alternative positions, or between  
           substituted amino acids. The nature of the cyclic  
           linkage may also be varied.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-14) - (G)4 - SDF-1(55-67) - K20/D24-cyclic acid

<400> 22

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15

Gly Gly Leu Lys Trp Ile Gln Asp Tyr Leu Glu Lys Ala Leu Asn  
20 25 30

<210> 23

<211> 31

<212> PRT

<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (15)..(18)

<223> spacer monomers (such as the illustrated glycine  
G's) may be used in variable numbers, such as 2, 3  
or 4 glycines.

<220>

<221> DOMAIN

<222> (20)..(24)

<223> Internal cyclization of peptides of the invention  
may be in alternative positions, or between  
substituted amino acids. The nature of the cyclic  
linkage may also be varied.

<220>

<223> Synthesised in Laboratory:

SDF-1(1-14) - (G)4 - SDF-1(55-67) - K20/D24-cyclic amide

<220>

<221> MOD\_RES

<222> (31)

<223> AMIDATION

<400> 23

Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15

Gly Gly Leu Lys Trp Ile Gln Asp Tyr Leu Glu Lys Ala Leu Asn  
20 25 30

<210> 24

<211> 31

<212> PRT

<213> Artificial Sequence

<220>

<221> DOMAIN

<222> (15)..(18)

<223> spacer monomers (such as the illustrated glycine  
G's) may be used in variable numbers, such as 2, 3  
or 4 glycines.

<220>  
 <221> DISULFID  
 <222> (9)..(11)  
 <223> cystein residues may for example be involved in  
 bridge formation

<220>  
 <223> Synthesised in Laboratory:  
 SDF-1(1-14) - (G)4-SDF-1(55-67) -C9/C11-cyclic acid

<400> 24  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
 1 5 10 15  
 Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
 20 25 30

<210> 25  
 <211> 31  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> DOMAIN  
 <222> (15)..(18)  
 <223> spacer monomers (such as the illustrated glycine  
 G's) may be used in variable numbers, such as 2, 3  
 or 4 glycines.

<220>  
 <221> DISULFID  
 <222> (9)..(11)  
 <223> Cysteine residues may for example be invloved in  
 bridge formation.

<220>  
 <223> Synthesised in Laboratory:  
 SDF-1(1-14) - (G)4-SDF-1(55-67) -C9/C11-cyclic amide

<220>  
 <221> MOD\_RES  
 <222> (31)  
 <223> AMIDATION

<400> 25  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
 1 5 10 15  
 Gly Gly Leu Lys Trp Ile Gln Glu Tyr Leu Glu Lys Ala Leu Asn  
 20 25 30

<210> 26  
 <211> 33  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <223> Synthesised in Laboratory: SDF-1(1-14) - (G)4-MIP-1  
 alpha(36-50)acid or amide.

<220>  
<221> MOD\_RES  
<222> (33)  
<223> Possible Amidation

<400> 26  
Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15  
Gly Gly Ser Lys Pro Gly Val Ile Phe Leu Thr Lys Arg Ser Arg Gln  
20 25 30

Val

<210> 27  
<211> 58  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory: SDF-1(1-14) - (G)4-MIP-1  
alpha(11-50)-acid or amide

<220>  
<221> MOD\_RES  
<222> (58)  
<223> Possible Amidation

<400> 27  
Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
1 5 10 15  
Gly Gly Cys Cys Phe Ser Tyr Thr Ser Arg Gln Ile Pro Gln Asn Phe  
20 25 30  
Ile Ala Asp Tyr Phe Glu Thr Ser Ser Gln Cys Ser Lys Pro Gly Val  
35 40 45  
Ile Phe Leu Thr Lys Arg Ser Arg Gln Val  
50 55

<210> 28  
<211> 33  
<212> PRT  
<213> Artificial Sequence

<220>  
<223> Synthesised in Laboratory: SDF-1(1-14) - (G)4-MIP-1  
alpha(56-70)-acid or amide

<220>  
<221> MOD\_RES  
<222> (33)  
<223> Possible Amidation

<400> 28  
 Lys Pro Val Ser Leu Ser Tyr Arg Cys Pro Cys Arg Phe Phe Gly Gly  
           1                  5                  10                  15

Gly Gly Glu Glu Trp Val Gln Lys Tyr Val Asp Asp Leu Glu Leu Ser  
                   20                  25                  30

Ala

<210> 29  
 <211> 9  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <221> BINDING  
 <222> (9)  
 <223> Lysine bridge may or may not be present between each of two arg  
           residues at position 8 of each of two monomers thereby forming a dimer.

<220>  
 <223> Synthesised in Laboratory: SDF-1(1-8)2-lysine  
           bridge dimer: or CTCE9904

<220>  
 <221> MOD\_RES  
 <222> (9)  
 <223> AMIDATION

<400> 29  
 Lys Pro Val Ser Leu Ser Tyr Arg Lys  
           1                  5

<210> 30  
 <211> 40  
 <212> PRT  
 <213> Artificial Sequence

<220>  
 <223> Synthesised in Laboratory

<400> 30  
 Cys Cys Phe Ser Tyr Thr Ser Arg Gln Ile Pro Gln Asn Phe Ile Ala Asp Tyr Phe  
           1                  5                  10                  15

Glu Thr Ser Ser Gln Cys Ser Lys Pro Gly Val Ile Phe Leu Thr Lys Arg  
           20                  25                  30                  35

Ser Arg Gln Val  
                   40

<210> 31  
 <211> 33  
 <212> PRT  
 <213> Artificial Sequence

<220>

<223> Synthesised in Laboratory: SDF-1(1-14) - (G)4-MIP-1  
alpha(36-50)-acid

<400> 31

Lys	Pro	Val	Ser	Leu	Ser	Tyr	Arg	Cys	Pro	Cys	Arg	Phe	Phe	Gly	Gly
1				5					10					15	

Gly	Gly	Ser	Lys	Pro	Gly	Val	Ile	Phe	Leu	Thr	Lys	Arg	Ser	Arg	Gln
			20					25					30		

Val